

Inhomogeneous structures of neutron star crust and mechanical properties

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Matter of neutron stars

Density

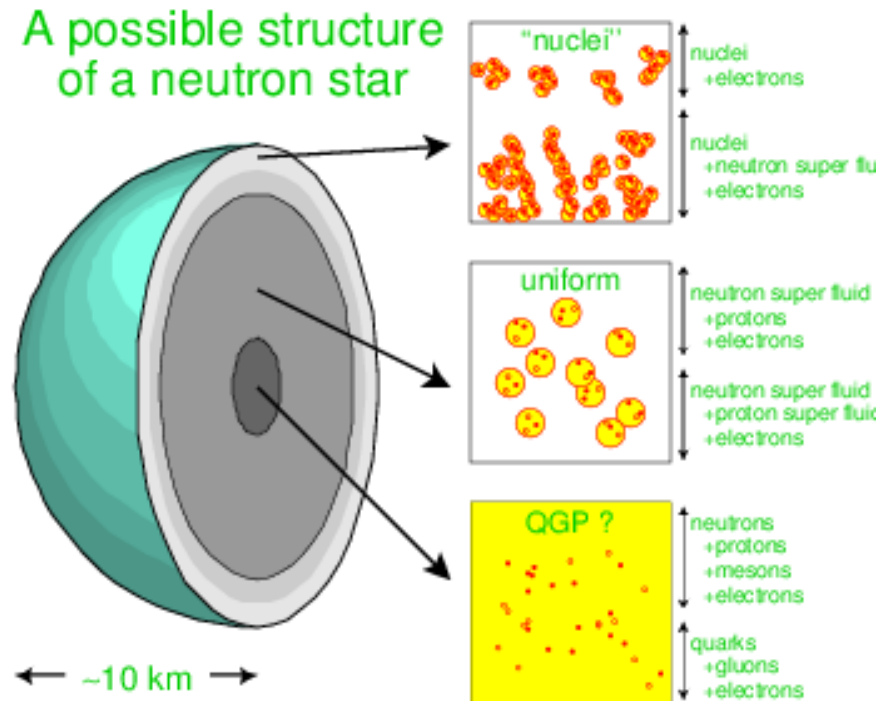
$$\approx 0 \dots \approx 10\rho_0$$

Composition

- Nucleons + leptons
- ... + mesons, hyperons
- quarks + gluons + leptons

Structure & correlation

- uniform
- crystal
- pasta
- amorphous
- pairing
-



Properties of neutron star matter

Compressibility (Equation of state).

- Masses and radii of neutron stars
- ← Hadron-quark phase transition
- ← Hyperon mixture and/or meson condensation
- ← Effects of mixed phase

Rigidity (shear modulus)

- Torsional oscillation might be directly measured as QPO
- Mountain height
- Starquakes

Our mean field studies of nuclear matter

- (inhomogeneous) structures of nuclear matter
- Eos
- **Useful for studies of mechanical strength (rigidity)**

RMF + Thomas-Fermi model

Lagrangian

$$L = L_N + L_M + L_e,$$

$$L_N = \bar{\Psi} \left[i\gamma^\mu \partial_\mu - m_N^* - g_{\omega N} \gamma^\mu \omega_\mu - g_{\rho N} \gamma^\mu \vec{\tau} \vec{b}_\mu - e \frac{1+\tau_3}{2} \gamma^\mu V_\mu \right] \Psi$$

$$L_M = \frac{1}{2} (\partial_\mu \sigma)^2 - \frac{1}{2} m_\sigma^2 \sigma^2 - U(\sigma) - \frac{1}{4} \omega_{\mu\nu} \omega^{\mu\nu} + \frac{1}{2} m_\omega^2 \omega_\mu \omega^\mu - \frac{1}{4} \vec{R}_{\mu\nu} \vec{R}^{\mu\nu} + \frac{1}{2} m_\rho^2 \vec{R}_\mu \vec{R}^\mu,$$

$$L_e = -\frac{1}{4} V_{\mu\nu} V^{\mu\nu} + \bar{\Psi}_e \left[i\gamma^\mu \partial_\mu - m_e + e\gamma^\mu V_\mu \right] \Psi_e, \quad (F_{\mu\nu} \equiv \partial_\mu F_\nu - \partial_\nu F_\mu)$$

$$m_N^* = m_N - g_{\sigma N} \sigma, \quad U(\sigma) = \frac{1}{3} b m_N (g_{\sigma N} \sigma)^3 + \frac{1}{4} c (g_{\sigma N} \sigma)^4$$

Nucleons interact with each other via coupling with σ , ω , ρ mesons.
Simple but feasible!

From $\partial_\mu \left[\partial L / \partial (\partial_\mu \phi) \right] - \partial L / \partial \phi = 0,$

$$(\phi = \sigma, \omega_\mu, R_\mu, V_\mu, \Psi),$$

$$-\nabla^2 \sigma(\mathbf{r}) + m_\sigma^2 \sigma(\mathbf{r}) = g_{\sigma N} (\rho_n^{(s)}(\mathbf{r}) + \rho_p^{(s)}(\mathbf{r})) - \frac{dU}{d\sigma}(\mathbf{r}),$$

$$-\nabla^2 \omega_0(\mathbf{r}) + m_\omega^2 \omega_0(\mathbf{r}) = g_{\omega N} (\rho_p(\mathbf{r}) + \rho_n(\mathbf{r})),$$

$$-\nabla^2 R_0(\mathbf{r}) + m_\rho^2 R_0(\mathbf{r}) = g_{\rho N} (\rho_p(\mathbf{r}) - \rho_n(\mathbf{r})),$$

$$\nabla^2 V_C(\mathbf{r}) = 4\pi e^2 \rho_{\text{ch}}(\mathbf{r}),$$

For Fermions, we employ Thomas-Fermi approx. with finite T

$$f_{i=n,p}(\mathbf{r}; \mathbf{p}, \mu_i) = \left(1 + \exp \left[\left(\sqrt{p^2 + m_i^*(\mathbf{r})^2} - \sqrt{p_{Fi}(\mathbf{r})^2 + m_i^*(\mathbf{r})^2} \right) / T \right] \right)^{-1},$$

$$f_e(\mathbf{r}; \mathbf{p}, \mu_e) = \left(1 + \exp \left[(p - (\mu_e - V_C(\mathbf{r}))) / T \right] \right)^{-1},$$

$$\rho_{i=p,n,e,v}(\mathbf{r}) = 2 \int_0^\infty \frac{d^3 p}{(2\pi)^3} f_i(\mathbf{r}; \mathbf{p}, \mu_i),$$

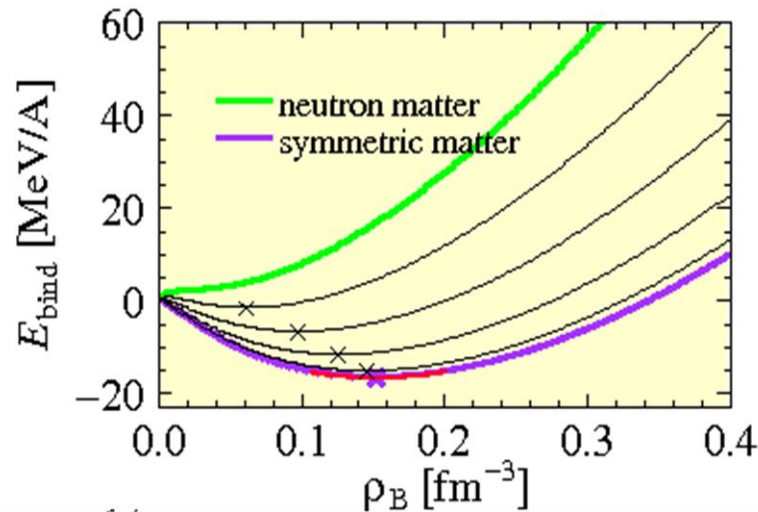
$$\mu_n = \sqrt{p_{Fn}(\mathbf{r})^2 + m_N^*(\mathbf{r})^2} + g_{\omega N} \omega_0(\mathbf{r}) - g_{\rho N} R_0(\mathbf{r}), \quad \mu_n = \mu_p + \mu_e,$$

$$\mu_p = \sqrt{p_{Fp}(\mathbf{r})^2 + m_N^*(\mathbf{r})^2} + g_{\omega N} \omega_0(\mathbf{r}) + g_{\rho N} R_0(\mathbf{r}) - V_C(\mathbf{r}),$$

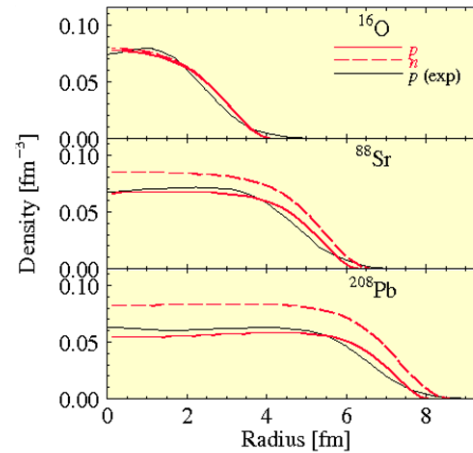
$$\int_V d^3 r \left[\rho_p(\mathbf{r}) + \rho_n(\mathbf{r}) \right] = \text{const}, \quad \int_V d^3 r \rho_p(\mathbf{r}) = \int_V d^3 r \rho_e(\mathbf{r}),$$

RMF + Thomas-Fermi model

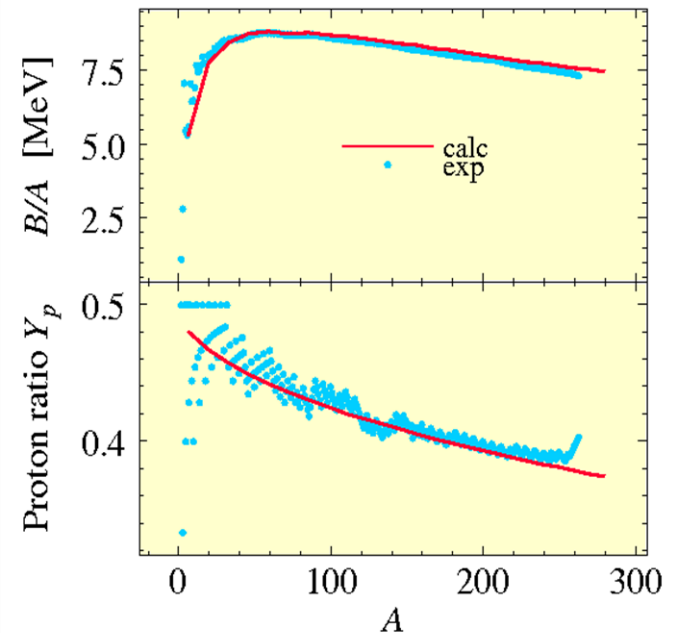
Nucleons interact with each other via coupling with σ , ω , ρ mesons. Simple but realistic enough.



Saturation property of symmetric nuclear matter : minimum energy $E/A \approx -16 \text{ MeV}$ at $\rho_B \approx 0.16 \text{ fm}^{-3}$.



Binding energies, proton fractions, and density profiles of nuclei are well reproduced.



Numerical procedure

To **equilibrate** $\mu_i(r)$ **in** r and **among species** i

$$\left. \begin{array}{l} \text{remove } r\text{-dependence } \mu_i(r) = \mu_i \\ \text{satisfy chemical balances } \mu_n = \mu_p + \mu_e \end{array} \right\} \#$$

- Divide whole space into equivalent and neutral cubic cells with **periodic boundary conditions**



- Distribute fermions (p, n, e) **randomly** but $\int d^3r \rho_i(r) = \text{given}$



- Solve field equations for $\sigma(r), \omega_0(r), \rho_0(r), V_{\text{Coul}}(r)$



- Calculate local chemical potentials of fermions $\mu_i(r)$



$$\mu_i(r) = V_i(r) + \sqrt{m_i^2 + p_{F_i}(r)^2}$$

- Adjust densities $\rho_i(r)$ as

$$\mu_i(r_1) > \mu_i(r_2) \quad \rightarrow \quad \rho_i(r_1) \downarrow, \rho_i(r_2) \uparrow$$

$$\mu_n(r) > \mu_p(r) + \mu_e \quad \rightarrow \quad \rho_n(r) \downarrow, \rho_p(r) \uparrow$$



- repeat until #

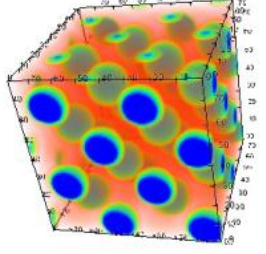
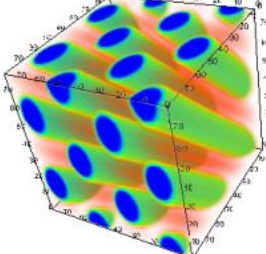
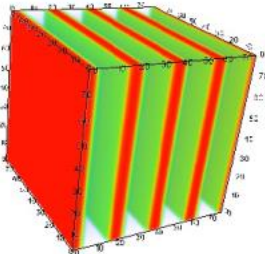
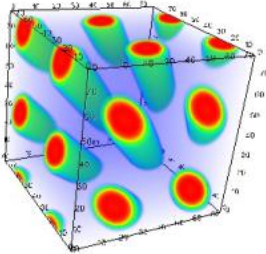
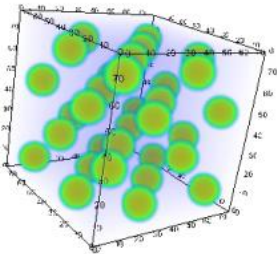
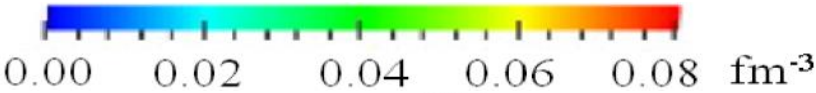


Fully 3D RMF calculations

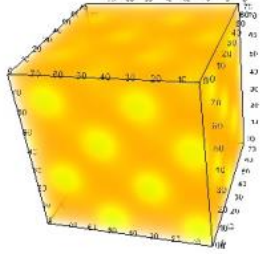
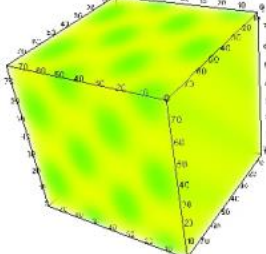
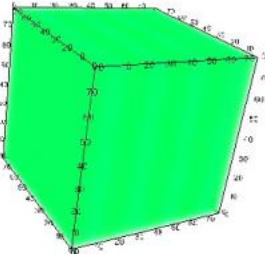
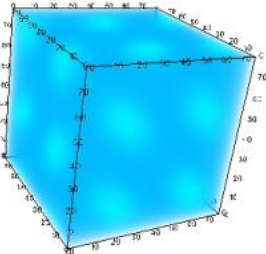
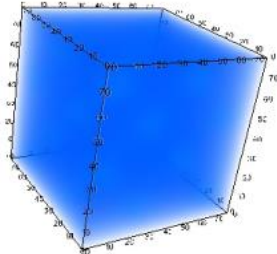
[Phys.Lett. B713 (2012) 284]

$$Y_p = Z/A = 0.5$$

proton



electron



“droplet”
[fcc]
 $\rho_B = 0.012 \text{ fm}^{-3}$

“rod”
 [honeycomb]
 0.024 fm^{-3}

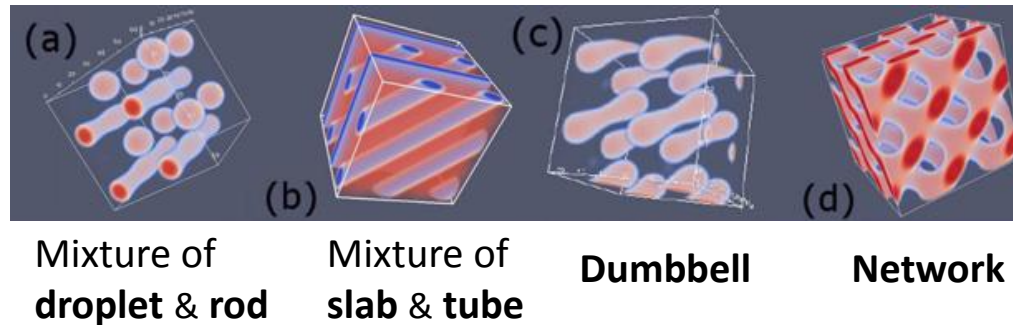
“slab”
 0.05 fm^{-3}

“tube”
 [honeycomb]
 0.08 fm^{-3}

“bubble”
 [fcc]
 0.094 fm^{-3}

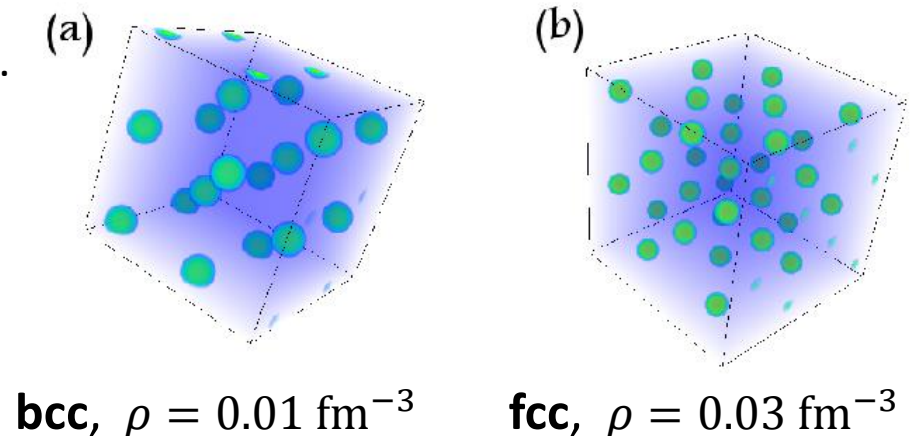
Advantages of a three-dimensional calculation

- Not only simple structures but any complex ones are taken into consideration in our new calculation.



Above structures are observed as excited states in our RMF calculation for symmetric ($Y_p = 0.5$) nuclear matter. Typical pasta structures are found to be the ground states.

- Crystalline structures can be discussed. We have found that fcc lattice appears at higher densities of droplet phase.



Shear

[Ogata et al,
PRA42(1990)4867]

$$r_i \rightarrow r_k \delta_{ik} u$$

$$(\vec{r} \rightarrow \vec{r}u)$$

Deformation = translation or rotation + isotropic compression + **shear**

Nothing essential

Conserves the volume.
There are 6 kinds

6 kinds of shear deformations

D1: $u_{xx} = \frac{2\epsilon - \epsilon^2}{(1-\epsilon)^2}, u_{yy} = u_{zz} = -\epsilon$

D2: D1 with $x \rightarrow y, y \rightarrow z, z \rightarrow x$

D3: D1 with $x \rightarrow z, y \rightarrow x, z \rightarrow y$

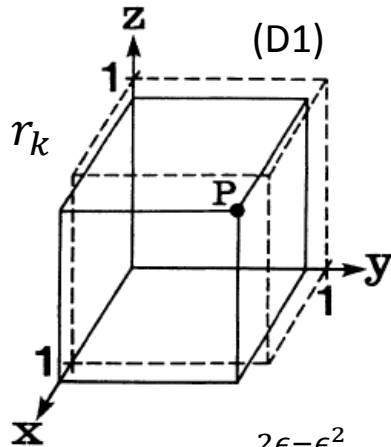
D4: $u_{xx} = \frac{\epsilon^2}{(1-\epsilon)^2}, u_{yz} = u_{zy} = \epsilon$

D5: D4 with $x \rightarrow y, y \rightarrow z, z \rightarrow x$

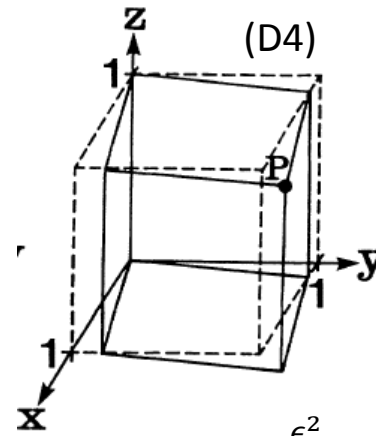
D6: D4 with $x \rightarrow z, y \rightarrow x, z \rightarrow y$

ϵ : infinitesimal

$$r_i \rightarrow (\delta_{ik} + u_{ik}) r_k$$



$$P = \left(1 + \frac{2\epsilon - \epsilon^2}{(1-\epsilon)^2}, 1 - \epsilon, 1 - \epsilon\right)$$



$$P = \left(1 + \frac{\epsilon^2}{1-\epsilon^2}, 1 + \epsilon, 1 + \epsilon\right)$$

Calculation of shear modulus

How rigid against shear deformation

increment of free energy

$$\delta F = \frac{1}{2} \lambda u_{ii}^2 + \mu u_{ik} u_{ik}$$

λ : volume elasticity

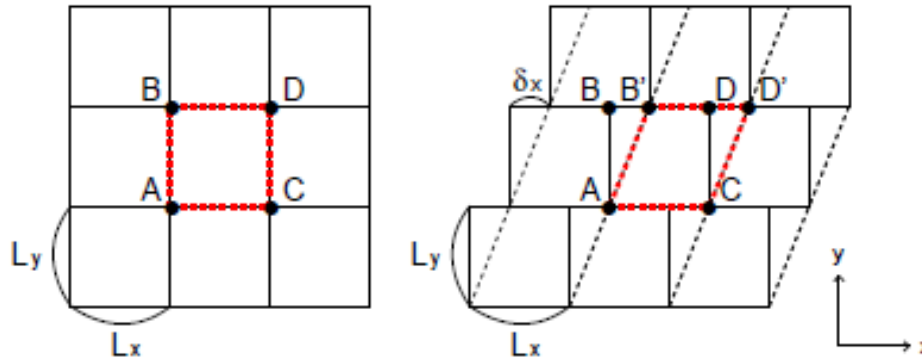
μ : shear modulus

How to calculate the shear modulus in our framework

- Prepare a ground state
- Give a small shear deformation without compression
- Calculate the curvature of the energy change against a shear deformation.
- Average the curvature among different shears

How to give a shear deformation

Deformed periodic boundary condition (DPBC)



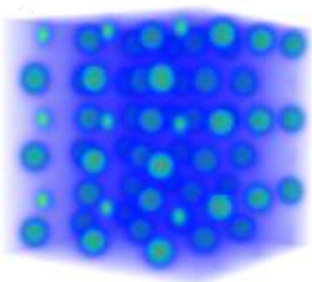
$$\epsilon = \frac{\delta x}{L_x}$$

$$F(x \pm L_x, y) = F(x, y), \quad F(x \pm L_x, y) = F(x, y),$$

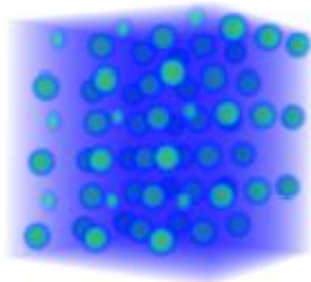
$$F(x, y \pm L_y) = F(x, y), \quad F(x, y \pm L_y) = F(x \mp \delta x, y)$$

Ground state under DPBC

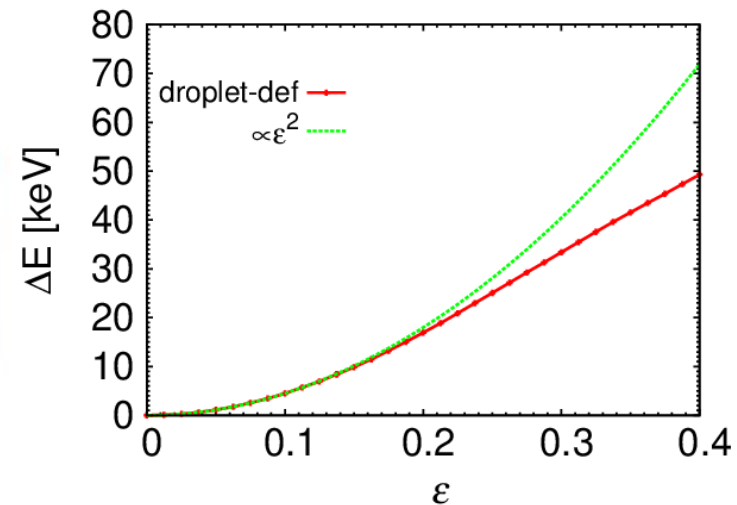
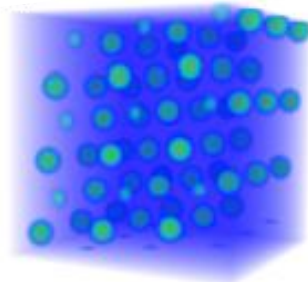
$\epsilon = 0$



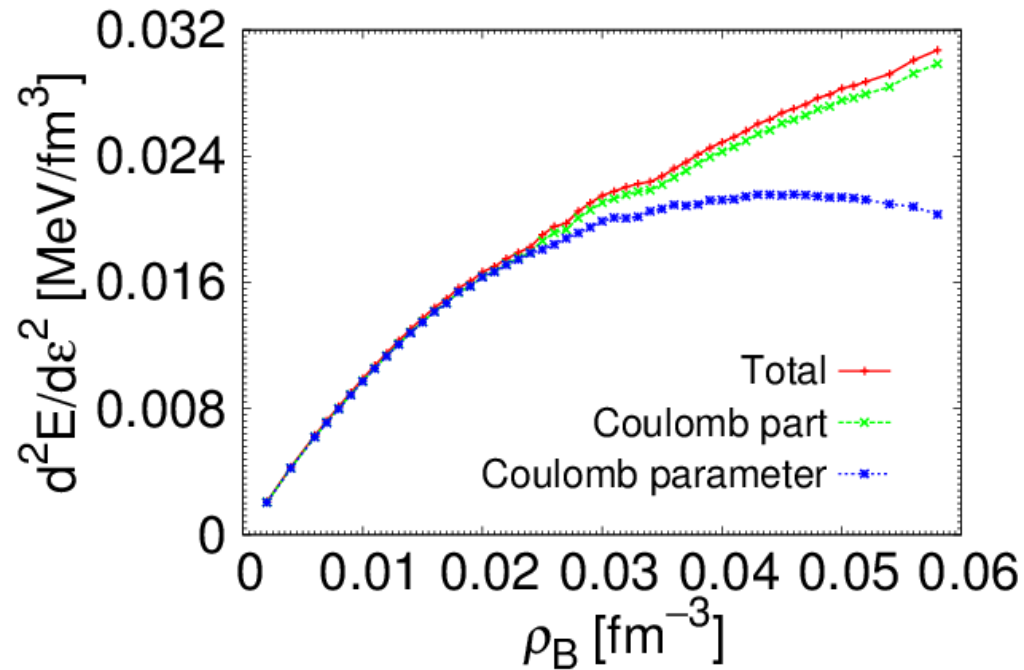
$\epsilon = 0.05$



$\epsilon = 0.1$



Preliminary result for droplet



- Below 0.02 fm^{-3} , we get shear modulus slightly smaller than the conventional study. \rightarrow screening effects.
- Above 0.02 fm^{-3} , our value is larger. \rightarrow finite size effects.

Summary

RMF + Thomas-Fermi calculation for neutron-star matter
Inhomogeneous structures appear below saturation density

Shear modulus for droplets calculated by deformed periodic boundary conditions
→ screening effects and finite-size effects

Future plan:

Comparison of shear moduli for bcc and fcc lattices and rod

